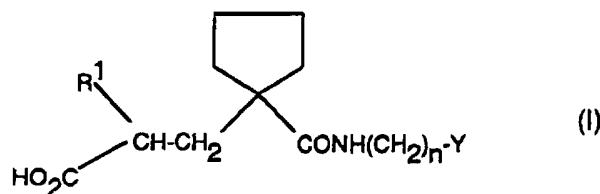


IN THE CLAIMS

Claim 1 (currently amended) A method of treating female sexual dysfunction comprising administering a therapeutically effective amount of a compound of formula (I), pharmaceutically acceptable salts, solvates, polymorphs or prodrugs thereof:



wherein

R^1 is C_{1-6} alkyl which may be substituted by one or more substituents, which may be the same or different, selected from the list: halo, hydroxy, C_{1-6} alkoxy, C_{2-6} hydroxyalkoxy, C_{1-6} alkoxy(C_{1-6} alkoxy), C_{3-7} cycloalkyl, C_{3-7} cycloalkenyl, aryl, aryloxy, (C_{1-4} alkoxy)aryloxy, heterocyclyl, heterocyclyloxy, $-NR^2R^3$, $-NR^4COR^5$, $-NR^4SO_2R^5$, $-CONR^2R^3$, $-S(O)_pR^6$, $-COR^7$ and $-CO_2(C_{1-4}$ alkyl); or R^1 is C_{3-7} cycloalkyl, aryl or heterocyclyl, each of which may be substituted by one or more substituents from said list, which substituents may be the same or different, which list further includes C_{1-6} alkyl; or R^1 is C_{1-6} alkoxy, $-NR^2R^3$ or $-NR^4SO_2R^5$;

wherein

R^2 and R^3 are each independently H, C_{1-4} alkyl, C_{3-7} cycloalkyl (optionally substituted by hydroxy or C_{1-4} alkoxy), aryl, (C_{1-4} alkyl)aryl, C_{1-6} alkoxyaryl or heterocyclyl; or R^2 and R^3 together with the nitrogen to which they are attached form a pyrrolidinyl, piperidino, morpholino, piperazinyl or N -(C_{1-4} alkyl)piperazinyl group;

R^4 is H or C_{1-4} alkyl;

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R^5 is C_{1-4} alkyl, CF_3 , aryl, $(C_{1-4}$ alkyl)aryl, $(C_{1-4}$ alkoxy)aryl, heterocyclyl,
 C_{1-4} alkoxy or $-NR^2R^3$ wherein R^2 and R^3 are as previously defined;

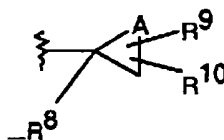
R^6 is C_{1-4} alkyl, aryl, heterocyclyl or NR^2R^3 wherein R^2 and R^3 are as
previously defined; and

R^7 is C_{1-4} alkyl, C_{3-7} cycloalkyl, aryl or heterocyclyl; p is 0, 1, 2 or 3;

n is 0, 1 or 2;

the $-(CH_2)_n-$ linkage is optionally substituted by C_{1-4} alkyl, C_{1-4} alkyl
substituted with one or more fluoro groups or phenyl, C_{1-4} alkoxy,
hydroxy, hydroxy(C_{1-3} alkyl), C_{3-7} cycloalkyl, aryl or heterocyclyl;

Y is the group

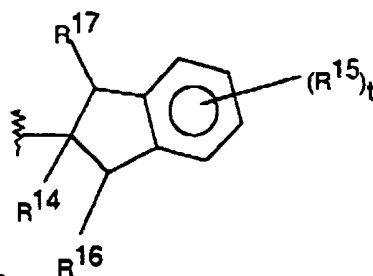


wherein A is $(CH_2)_q$ where q is 1, 2, 3 or 4 to complete a 3 to 7 membered
carbocyclic ring which may be saturated or unsaturated; R^8 is H, C_{1-6}
alkyl, CH_2OH , phenyl, phenyl(C_{1-4} alkyl) or $CONR^{11}R^{12}$; R^9 and
 R^{10} are each independently H, CH_2OH , $C(O)NR^{11}R^{12}$, C_{1-6} alkyl,
phenyl (optionally substituted by C_{1-4} alkyl, halo or C_{1-4} alkoxy or
phenyl(C_{1-4} alkyl) wherein the phenyl group is optionally substituted by
 C_{1-4} alkyl, halo or C_{1-4} alkoxy, or R^9 and R^{10} together form a
dioxolane; R^{11} and R^{12} which may be the same or different are H, C_{1-4}
alkyl, R^{13} or $S(O)_rR^{13}$, where r is 0, 1 or 2 and R^{13} is phenyl
optionally substituted by C_{1-4} alkyl or phenyl(C_{1-4} alkyl) wherein the
phenyl is optionally substituted by C_{1-4} alkyl; or

Y is the group, $C(O)NR^{11}R^{12}$ wherein R^{11} and R^{12} are as previously
defined except that R^{11} and R^{12} are not both H; or

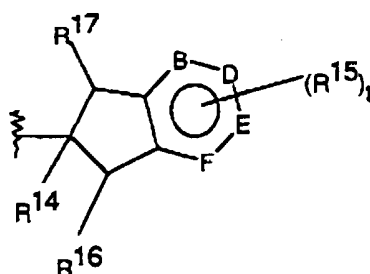
Y is the group,

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wherein R^{14} is H, CH_2OH , or $C(O)NR^{11}R^{12}$ wherein R^{11} and R^{12} are as previously defined; when present R^{15} , which may be the same or different to any other R^{15} , is OH, C_{1-4} alkyl, C_{1-4} alkoxy, halo or CF_3 ; t is 0, 1, 2, 3 or 4; and R^{16} and R^{17} are independently H or C_{1-4} alkyl;
or

Y is the group



wherein one or two of B, D, E or F is a nitrogen, the others being carbon; and
 R^{14} to R^{17} and t are as previously defined; or

Y is an optionally substituted 5-7 membered heterocyclic ring, which may be saturated, unsaturated or aromatic and contains a nitrogen, oxygen or sulphur and optionally one, two or three further nitrogen atoms in the ring and which may be optionally benzofused and optionally substituted by:

C_{1-6} alkoxy; hydroxy; oxo; amino; mono or di- $(C_{1-4}$ alkyl)amino;

C_{1-4} alkanoylamino; or

C_{1-6} alkyl which may be substituted by one or more substituents, which may be the same or different, selected from the list: C_{1-6} alkoxy, C_{1-6} haloalkoxy, C_{1-6} alkylthio, halogen, C_{3-7} cycloalkyl, heterocyclyl or phenyl; or

C₃₋₇cycloalkyl, aryl or heterocyclyl, each of which may be substituted by one or more substituents, which may be the same or different, selected from the list: C₁₋₆alkyl, C₁₋₆alkoxy, C₁₋₆haloalkoxy, C₁₋₆alkylthio, halogen, C₃₋₇cycloalkyl, heterocyclyl or phenyl;

wherein when there is an oxo substitution on the heterocyclic ring, the ring only contains one or two nitrogen atoms and the oxo substitution is adjacent a nitrogen atom in the ring; or

~~Y is $\text{NR}^{18}\text{S}(\text{O})_u\text{R}^{19}$, wherein R^{18} is H or C₁₋₄alkyl; R^{19} is aryl, arylC₁₋₄alkyl or heterocyclyl; and u is 0, 1, 2 or 3.~~

Claim 2 (currently amended) A compound of formula (I), pharmaceutically acceptable salts, solvates, polymorphs or prodrugs thereof, wherein R¹, n and Y are as defined in claim 1 with the proviso that ~~Y is not the group $\text{C}(\text{O})\text{NR}^{11}\text{R}^{12}$ and~~ when R¹ is propyl or phenylethyl, R¹⁴ is not -CH₂OH.

Claim 3 (currently amended) A compound of formula (I), pharmaceutically acceptable salts, solvates, polymorphs or prodrugs thereof, wherein R¹, n and Y are as defined in claim 1 with the proviso that ~~Y is not the group $\text{C}(\text{O})\text{NR}^{11}\text{R}^{12}$ and~~ R¹⁴ is not H or -CH₂OH.

Claim 4 (original) A compound according to claim 2, pharmaceutically acceptable salts, solvates, polymorphs or prodrugs thereof, wherein R¹ is C₁₋₆alkyl, C₁₋₆alkoxy, C₁₋₆alkoxy(C₁₋₃)alkyl, C₁₋₆alkoxyC₁₋₆alkoxyC₁₋₃alkyl or C₁₋₆alkyl substituted with aryl.

Claim 5 (original) A compound according to claim 4, pharmaceutically acceptable salts, solvates, polymorphs or prodrugs thereof, wherein R¹ is C₁₋₆alkyl, C₁₋₆alkoxy, C₁₋₆alkoxy(C₁₋₃)alkyl or C₁₋₆alkoxyC₁₋₆alkoxyC₁₋₃alkyl.

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Claim 6 (original) A compound according to claim 5, pharmaceutically acceptable salts, solvates, polymorphs or prodrugs thereof, wherein R¹ is C₁₋₄alkyl or C₁₋₆alkoxy(C₁₋₃)alkyl.

Claims 7-13 (withdrawn)

Claim 14 (Original) A compound according to claim 2, pharmaceutically acceptable salts, solvates, polymorphs or prodrugs thereof, wherein Y is an optionally substituted 5-7 membered heterocyclic ring.

Claim 15 (original) A compound according to claim 14, pharmaceutically acceptable salts, solvates, polymorphs or prodrugs thereof, wherein the 5-7 membered heterocyclic ring is an optionally substituted aromatic ring.

Claim 16 (original) A compound according to claim 15, pharmaceutically acceptable salts, solvates, polymorphs or prodrugs thereof, wherein said aromatic ring is pyridyl, pyrazinyl, pyrimidinyl, pyridazinyl, pyrazolyl, triazolyl, tetrazolyl, oxadiazolyl, thiazolyl, thiadiazolyl, oxazolyl, isoxazolyl, indolyl, isoindolyl, quinolyl, isoquinolyl, pyridonyl, quinoxalinyll or quinazolinyl each of which may be substituted as defined in claim 1.

Claim 17 (original) A compound according to claim 16, pharmaceutically acceptable salts, solvates, polymorphs or prodrugs thereof, wherein the aromatic ring is oxadiazole, pyridone or thiadiazole each of which may be substituted as defined in claim 1.

Claim 18 (original) A compound according to claim 17, pharmaceutically acceptable salts, solvates, polymorphs or prodrugs thereof, wherein the aromatic ring is 1,2,5-oxadiazole, 1,3,4-oxadiazole, 2-pyridone or 1,3,4-thiadiazole each of which may be substituted as defined in claim 1.

Claim 19 (original) A compound according to claim 14, pharmaceutically

acceptable salts, solvates, polymorphs or prodrugs thereof, wherein the 5-7 membered heterocyclic ring is substituted by one or more C₁₋₆alkyl, phenyl or phenylC₁₋₄alkyl.

Claim 20 (original) A compound according to claim 19, pharmaceutically acceptable salts, solvates, polymorphs or prodrugs thereof, wherein the 5-7 membered heterocyclic ring is substituted by C₁₋₄alkyl or benzyl.

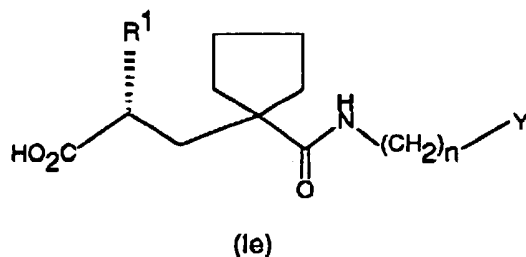
Claim 21 (original) A compound according to claim 17, pharmaceutically acceptable salts, solvates, polymorphs or prodrugs thereof, wherein when Y is a pyridone said pyridone is *N*-substituted pyridone.

Claims 22-23 (withdrawn)

Claim 24 (original) A compound according to claim 2, pharmaceutically acceptable salts, solvates, polymorphs or prodrugs thereof, wherein R¹⁶ and R¹⁷ are hydrogen.

Claim 25 (original) A compound according to claim 2, pharmaceutically acceptable salts, solvates, polymorphs or prodrugs thereof, wherein t is 0.

Claim 26 (original) A compound of formula 1e, pharmaceutically acceptable salts, solvates, polymorphs or prodrugs thereof,



wherein R¹, Y and n are as defined in claim 2.

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Claim 27 (currently amended) A compound, pharmaceutically acceptable salts, solvates, polymorphs or prodrugs thereof, selected from the group consisting of:

2-[(1-[(1-benzyl-6-oxo-1,6-dihydro-3-pyridinyl)amino]carbonyl)cyclopentyl]-methyl]-4-methoxybutanoic acid;

2-[[1-[(3-(2-oxo-1-pyrrolidinyl)propyl)amino]carbonyl(cyclopentyl)]-methyl]-4-phenylbutanoic acid);

(+)-2-[[1-[(2-(hydroxymethyl)-2,3-dihydro-1H-inden-2-yl)amino]carbonyl)cyclopentyl]methyl]-4-phenylbutanoic acid;

2-[(1-[(5-methyl-1,3,4-thiadiazol-2-yl)amino]carbonyl)cyclopentyl]methyl]-4-phenylbutanoic acid;

~~cis-3-(2-methoxyethoxy)-2-[(1-[(4-((phenylsulfonyl)amino)carbonyl)cyclohexyl]-amino]carbonyl)cyclopentyl]methyl]propanoic acid;~~

(+)-2-[[1-[(2-(hydroxymethyl)-2,3-dihydro-1H-inden-2-yl)amino]carbonyl)cyclopentyl]-methyl]pentanoic acid;

(2R)-2-[(1-[(5-ethyl-1,3,4-thiadiazol-2-yl)amino]carbonyl)cyclopentyl]-methyl]pentanoic acid or (-)-2-[(1-[(5-ethyl-1,3,4-thiadiazol-2-yl)amino]carbonyl)cyclopentyl]-methyl]pentanoic acid;

(2S)-2-[(1-[(5-ethyl-1,3,4-thiadiazol-2-yl)amino]carbonyl)cyclopentyl]-methyl]pentanoic acid or (+)-2-[(1-[(5-ethyl-1,3,4-thiadiazol-2-yl)amino]carbonyl)cyclopentyl]-methyl]pentanoic acid ;

~~2-[(1-[(3-benzylaniline)carbonyl)cyclopentyl]methyl]pentanoic acid ;~~

2-[(1-[(1-benzyl-6-oxo-1,6-dihydro-3-pyridinyl)amino]carbonyl)cyclopentyl]-methyl]pentanoic acid ;

~~2-[[1-[(1R,3S,4R)-4-(aminocarbonyl)-3-butylcyclohexyl]amino]carbonyl]-cyclopentyl]methyl]pentanoic acid ;~~

~~trans-3-[1-[(2-(4-chlorophenyl)cyclopropyl)amino]carbonyl)cyclopentyl]-2-(methoxymethyl)propanoic acid ;~~

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~~trans-3-[1-(((2-(4-methoxyphenyl)cyclopropyl)amino)carbonyl)cyclopentyl]-2-(methoxyethyl)propanoic acid;~~
~~trans-3-[1-(((2-pentylcyclopropyl)amino)carbonyl)cyclopentyl]-2-(methoxyethyl)propanoic acid;~~
 3-[1-(((5-benzyl-[1,3,4]-thiadiazol-2-yl)amino)carbonyl)cyclopentyl]-2-(methoxyethyl)propanoic acid ;
 3-[1-(((4-butylpyridin-2-yl)amino)carbonyl)cyclopentyl]-2-(methoxyethyl)propanoic acid ;
 3-[1-(((4-phenylpyridin-2-yl)amino)carbonyl)cyclopentyl]-2-(methoxyethyl)propanoic acid ;
~~3-[1-(((1-hydroxymethyl-3-phenylcyclopentyl)amino)carbonyl)cyclopentyl]-2-(methoxyethyl)propanoic acid;~~
 2-[[1-(((2-(hydroxymethyl)-2,3-dihydro-1*H*-inden-2-yl)amino)carbonyl)-cyclopentyl]methyl]-4-methoxybutanoic acid ;
~~trans-3-[1-(((2-phenylcyclopropyl)amino)carbonyl)cyclopentyl]-2-(methoxyethyl)propanoic acid;~~
 (*R*)-2-[[1-(((2-(hydroxymethyl)-2,3-dihydro-1*H*-inden-2-yl)amino)carbonyl)-cyclopentyl]methyl]-4-methoxybutanoic acid ; and
 (*S*)-2-[[1-(((2-(hydroxymethyl)-2,3-dihydro-1*H*-inden-2-yl)amino)carbonyl)-cyclopentyl]methyl]-4-methoxybutanoic acid .

Claim 28 (original) The method according to claim 1 wherein the female sexual dysfunction treated includes at least female sexual arousal dysfunction (FSAD).

Claim 29 (original) The method according to claim 1 wherein the medicament is administered systemically.

Claim 30 (original) The method according to claim 1 wherein the medicament is administered orally.

Claim 31 (currently amended) A method of treatment or prophylaxis of a condition for which a beneficial therapeutic response can be obtained by the inhibition of

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neutral endopeptidase comprising administration of a therapeutically effective amount of a compound as defined in claim 2.

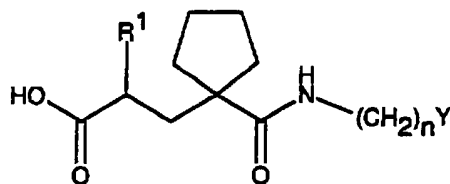
Claim 32. (Previously Cancelled)

Claim 33. (Previously Amended) A pharmaceutical formulation comprising a compound as defined in claim 2 together with a pharmaceutically acceptable excipient.

Claim 34. (Previously Amended) A method for the treatment or prophylaxis of female sexual dysfunction comprising administering to the patient a therapeutically effective amount of a compound as defined in claim 2.

Claim 35. (Previously Cancelled)

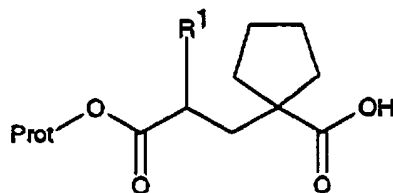
Claim 36 (currently amended) A process for preparing a compound of formula I or salts thereof



(I)

wherein R^1 , n and Y are as defined in any one of claims 2 to 27, comprising the steps of:

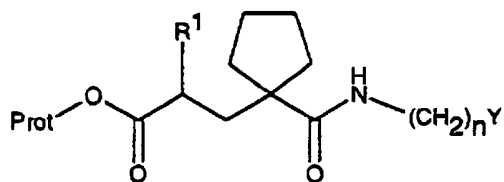
a) reacting a compound of formula II



(II)

wherein Prot is a suitable protecting group, with a compound of formula $\text{Y}(\text{CH}_2)_n\text{NH}_2$ (III), to give a compound of formula IV,

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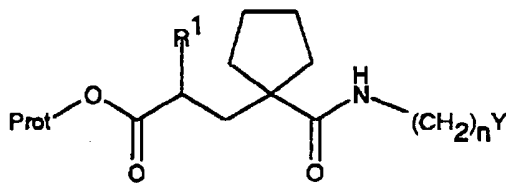


(IV)

then

- b) reacting the compound of formula IV under suitable deprotecting conditions to give the compound of formula I; then
- c) optionally forming a salt.
- d)

Claim 37. (original) A compound of formula IV



(IV)

wherein R^1 , n , and Y are as defined in claim 2 and wherein Prot is a protecting group.